This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1) (Previously presented) A compound of formula (I):

wherein

A is a bicyclic heterocycle which is:

- (1) benzimidazolyl
- (2) 1,3-benzothiazolyl
- (3) 1,2,3-benzotriazolyl
- (4) 1,3-benzoxazolyl
- (5) 2,3-dihydro-1H-indolyl
- (6) 2,3-dihydro-1H-indenyl
- (7) 1,1-dioxido-2,3-dihydro-1-benzothienyl
- (8) 1H-indazolyl
- (9) 2H-indazolyl
- (10) 1H-indolyl
- (11) 2H-chromenyl
- (12) quinoxalinyl or
- (13) a group of the formula

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optionally substituted with 1-4 substituents which are independently R^1 , OR^1 , $S(O)_pR^1$, $C(O)R^1$, $C(O)OR^1$, $C(O)NR^1R^2$, halogen, oxo, cyano, or nitro;

B is phenyl, naphthyl or, pyridyl, optionally substituted with 1-4 substituents which are independently C_1 - C_5 linear or branched alkyl, C_1 - C_5 linear or branched haloalkyl, C_1 - C_3 alkoxy, hydroxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, carboxyamide, halogen, cyano, nitro or $S(O)_0R^7$;

Lis:

- (a) -(CH₂)_m-O-(CH₂)_l-,
- (b) -(CH₂)_m-(CH₂)_l-,
- (c) -(CH₂)_m-C(O)-(CH₂)₁-.
- (d) -(CH₂)_m-NR³-(CH₂)₁-,
- (e) -(CH₂)_m- NR³C(O)-(CH₂)_i-,
- (f) -(CH₂)_m-S-(CH₂)_l-,
- (g) -(CH₂)_m-C(O)NR³ -(CH₂)_i-, or
- (h) a single bond:

m and I are integers independently selected from 0-4;

M is a pyridine ring, optionally substituted with 1-3 substituents which are independently C_1 - C_5 linear or branched alkyl, C_1 - C_5 linear or branched haloalkyl, C_1 - C_3 alkoxy, hydroxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, halogen, or nitro;

Q is C(O)R4, C(O)OR4 or C(O)NR4R5.;

each of R1, R2, R3, R4 and R5 is independently:

- (a) hydrogen,
- (b) C₁-C₅ linear, branched, or cyclic alkyl,
- (c) phenyl,

- (d) C₁-C₃ alkyl-phenyl,
- (e) up to per-halo substituted C₁-C₅ linear or branched alkyl,
- (f) -(CH₂)_q·X, wherein X is a 5 or 6 membered heterocyclic ring, containing at least one atom selected from oxygen, nitrogen and sulfur, which is saturated, partially saturated, or aromatic, or a 8-10 membered bicyclic heteroaryl having 1-4 heteroatoms which are O, N or S, or
- (g) -(CH₂)_g-Y, where Y is C(O)R⁶, C(O)OR⁶ and C(O)NR⁶R⁷;

each of R⁶ - R⁷ is independently:

- (a) hydrogen,
- (b) C₁-C₅ linear, branched, or cyclic alkyl,
- (c) phenyl,
- (d) C₁-C₃ alkyl-phenyl, or
- (e) up to per-halo substituted C₁-C₅ linear or branched alkyl;

each of R¹, R², R³, R⁴, R⁵, R⁶ and R⁷, other than per–halo substituted C_1 - C_5 linear or branched alkyl, is optionally substituted with 1-3 substituents which are independently C_1 - C_5 linear or branched alkyl, up to perhalo substituted C_1 - C_5 linear or branched alkyl, C_1 - C_3 alkoxy, hydroxy, carboxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, halogen, cyano, or nitro;

p is an integer selected from 0, 1, or 2; and

q is an integer selected from 1, 2, 3, or 4

or a pharmaceutically acceptable salt of formula I or an oxidized derivative of formula I wherein one or more urea nitrogens are substituted with a hydroxyl group, or an oxidized derivative of formula I wherein the nitrogen atom of pyridine ring M is in the oxide form, or a methyl, ethyl, propyl, isopropyl, butyl, isobutyl, pentyl ester or phenyl C₁-C₅ alkyl ester of formula I at a carboxylic acid group or amide group.

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 (previously presented) A compound of claim 1 wherein A and B follow one of the following combinations:

A= 1H-benzimidazol-5-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1H-benzimidazol-6-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,3-benzodioxin-6-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,3-benzodioxin-7-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,3-benzodioxin-8-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,3-benzodioxol-4-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,3-benzodioxol-5-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,3-benzothiazol-2-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,3-benzothiazol-5-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,3-benzothiazol-6-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,2,3-benzotriazol-5-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1,3-benzoxazol-2-yl; and B= phenyl, pyridinyl or naphthyl, or

A= 1,3-benzoxazol-6-yl; and B= phenyl, pyridinyl or naphthyl.

3) (original) A compound of claim 1 wherein A and B follow one of the following combinations:

A= 1H-benzimidazolyl; and B= phenyl or pyridinyl,

A= 1,3-benzodioxinyl; and B= phenyl or pyridinyl,

A= 1,3-benzodioxolyl; and B= phenyl or pyridinyl,

A= 1,3-benzothiazolyl; and B= phenyl or pyridinyl,

A= 1,2,3-benzotriazolyl; and B= phenyl or pyridinyl, or

A= 1,3-benzoxazolyl; and B= phenyl, pyridinyl.

 (original) A compound of claim 1 wherein A and B follow one of the following combinations:

A= 1H-benzimidazol-5-yl; and B= phenyl or pyridinyl,

A= 1H-benzimidazol-6-yl; and B= phenyl or pyridinyl,

A= 1,3-benzodioxin-6-yl; and B= phenyl or pyridinyl,,
A= 1,3-benzodioxin-7-yl; and B= phenyl or pyridinyl.

A= 1,3-benzodioxin-7-yi; and B= pnenyi or pyridiny

- A= 1,3-benzodioxin-8-yl; and B= phenyl or pyridinyl,
- A= 1,3-benzodioxol-4-yl; and B= phenyl or pyridinyl, ,
- A= 1,3-benzodioxol-5-yl; and B= phenyl or pyridinyl,
- A=1,3-benzothiazol-2-yl; and B= phenyl or pyridinyl,
- A=1,3-benzothiazol-5-yl; and B= phenyl or pyridinyl,
- A= 1,3-benzothiazol-6-yl; and B= phenyl or pyridinyl, A= 1,2,3-benzotriazol-5-yl; and B= phenyl or pyridinyl,
- A= 1,3-benzoxazol-2-yl; and B= phenyl or pyridinyl, or
- A= 1,3-berizoxazoi-z-yi, and b= prienyi or pyridinyi, or
- A= 1.3-benzoxazol-6-vl; and B= phenyl or pyridinyl.
- 5) (previously presented) A compound of claim 1 wherein A and B follow one of the following combinations:
- $A = 2, 3 dihydro-1, 4 benzodioxin-5 yl; \ and \ B = \ phenyl, \ pyridinyl \ or \\ naphthyl,$
- $A = 2, 3 dihydro 1, 4 benzodioxin 6 yl; \quad and \quad B = \quad phenyl, \quad pyridinyl \quad or \\ naphthyl,$
 - A= 2,3-dihydro-1-benzofuran-5-yl; and B= phenyl, pyridinyl or naphthyl,
 - A= 2,3-dihydro-1H-indol-5-yl; and B= phenyl, pyridinyl or naphthyl,
 - A=2,3-dihydro-1H-indol-6-yl; and B= phenyl, pyridinyl or naphthyl,
 - A=2,3-dihydro-1H-inden-4-yl; and B= phenyl, pyridinyl or naphthyl,
 - A= 2,3-dihydro-1H-inden-5-yl; and B= phenyl, pyridinyl or naphthyl,
- $A = 1,1-dioxido-2,3-dihydro-1-benzothien-6-yl; \ and \ B = phenyl, \ pyridinyl or naphthyl.$
- 6) (original) A compound of claim 1 wherein A and B follow one of the following combinations:
 - A= 2,3-dihydro-1,4-benzodioxin-5-yl; and B= phenyl or pyridinyl,
 - A= 2,3-dihydro-1,4-benzodioxin-6-yl; and B= phenyl or pyridinyl,
 - A= 2,3-dihydro-1-benzofuran-5-yl; and B= phenyl or pyridinyl, A= 2,3-dihydro-1H-indol-5-yl; and B= phenyl or pyridinyl,
 - A= 2,3-dihydro-1H-indol-6-yl; and B= phenyl or pyridinyl,
 - A= 2,3-dihydro-1H-inden-4-yl; and B= phenyl or pyridinyl,

A= 2,3-dihydro-1H-inden-5-yl; and B= phenyl or pyridinyl, or A= 1,1-dioxido-2,3-dihydro-1-benzothien-6-yl; and B= phenyl or pyridinyl.

7) (previously presented) A compound of claim 1 wherein A and B follow one of the following combinations:

A= 1H-indazol-5-yl; and B= phenyl, pyridinyl or naphthyl,

A= 2H-indazol-5-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1H-indazol-6-yl; and B= phenyl, pyridinyl or naphthyl,

A= 1H-indol-5-yl; and B= phenyl, pyridinyl or naphthyl,

A= 2-oxo-2H-chromen-7-yl; and B= phenyl, pyridinyl or naphthyl or

 $\mbox{A= 1-oxo-2,3-dihydro-1H-inden-5-yl; and B= phenyl, pyridinyl or} \label{eq:AB} \mbox{naphthyl.}$

8) (original) A compound of claim 1 wherein A and B follow one of the following combinations:

A= 1H-indazol-5-yl; and B= phenyl or pyridinyl,

A= 2H-indazol-5-yl; and B= phenyl or pyridinyl,

A= 1H-indazol-6-yl; and B= phenyl or pyridinyl,

A= 1H-indol-5-yl; and B= phenyl or pyridinyl,

A= 2-oxo-2H-chromen-7-yl; and B= phenyl or pyridinyl, or

A= 1-oxo-2,3-dihydro-1H-inden-5-yl; and B= phenyl or pyridinyl.

9) (previously presented) A compound of claim 1 wherein A and B follow one of the following combinations:

A= quinoxalin-2-yl; and B= phenyl, pyridinyl or naphthyl or A= quinoxalin-6-yl; and B= phenyl, pyridinyl or naphthyl.

10) (original) A compound of claim 1 wherein A and B follow one of the following combinations:

A= quinoxalin-2-yl; and B= phenyl or pyridinyl, or A= quinoxalin-6-yl; and B= phenyl or pyridinyl.

- 11) (original) A compound as in claim 1 wherein L is -O- or -S-.
- 12) (previously presented) A compound which is:
- N-methyl-4-[3-({[(2-methyl-1,3-benzoxazol-6yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- 4-[4-({[(1-acetyl-2,3-dihydro-1H-indol-6-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-([[(6-chloro-1,3-benzothiazol-2-yl)amino]carbonyl}amino)phenoxy]-Nmethylpyridine-2-carboxamide
- N-methyl-4-{4-[([6-(trifluoromethoxy)-1,3-benzothiazol-2yl]amino}carbonyl)amino]phen-oxy}pyridine-2-carboxamide
- 4-[4-([[(6-fluoro-1,3-benzothiazol-2-yl)amino]carbonyl]amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[3-fluoro-4-({[(6-fluoro-1,3-benzothiazol-2-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-{3-fluoro-4-[({{6-(trifluoromethoxy)-1,3-benzothiazol-2-yl]amino}carbonyl)amino]phen-oxy}-N-methylpyridine-2-carboxamide;
- 4-[4-({[(6-methoxy-1,3-benzothiazol-2-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-({[(6-methoxy-1,3-benzothiazol-2-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-([[(5-chloro-1,3-benzoxazol-2-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-([[(5-chloro-1,3-benzoxazol-2-yl)amino]carbonyl]amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-({[(6-chloro-1,3-benzothiazol-2-yl)amino]carbonyl}amino)-3fluorophenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-({[(6-chloro-1,3-benzothiazol-2-yl)amino]carbonyl]amino)-3fluorophenoxy]-N-methylpyridine-2-carboxamide

- 4-(2-chloro-4-[[(2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino)phenoxy)-N-methylpyridine-2-carboxamide
- 4-[(5-[((2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino)quinolin-8-yl)oxy]-N-methylpyridine-2-carboxamide
- 4-[4-({[(4,6-difluoro-1,3-benzothiazol-2-yl)amino]carbonyl}amino)-3fluorophenoxy]-N-methylpyridine-2-carboxamide
- 4-[3-fluoro-4-([(6-methoxy-1,3-benzothiazol-2yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-(4-{[({1-[2-(diethylamino)ethyl]-1H-indol-5-yl}amino)carbonyl]amino}-3fluorophenoxy)-N-methylpyridine-2-carboxamide;
- 4-(4-{[(2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino}-3-fluorophenoxy)-N-methylpyridine-2-carboxamide
- 4-[3-fluoro-4-([[(1-oxo-2,3-dihydro-1H-inden-5yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-([[(1,1-dioxido-2,3-dihydro-1-benzothien-6-yl)amino]carbonyl}amino)-3-fluorophenoxyl-N-methylpyridine-2-carboxamide
- 4-[3-fluoro-4-([[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[2-fluoro-4-([[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[2,4-difluoro-5-([[(1-methyl-1H-indazol-5yl)amino]carbonyl]amino)phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-[4-([[(1-methyl-1H-indazol-5-yl)amino]carbonyl]amino)-3-(trifluoromethyl)-phenoxy]pyridine-2-carboxamide
- 4-[4-fluoro-3-([[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[2-fluoro-5-([[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[2-chloro-6-fluoro-4-({[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[3-fluoro-4-([[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-(2-methoxyethyl)pyridine-2-carboxamide

- 4-[3-fluoro-4-({[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-({[(2,2-difluoro-1,3-benzodioxol-5-yl)amino]carbonyl}amino)-3fluorophenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-(4-{[(quinoxalin-6-ylamino)carbonyl]amino}phenoxy)pyridine-2carboxamide
- 4-(3-fluoro-4-{[(quinoxalin-6-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- N-methyl-4-[4-{[(quinoxalin-6-ylamino)carbonyl]amino}-3-(trifluoromethyl)phenoxyl-pyridine-2-carboxamide
- 4-(3-chloro-4-{[(quinoxalin-6-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- N-methyl-4-[4-({[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]carbonyl}-amino)-3-(trifluoromethyl)phenoxy]pyridine-2-carboxamide
- 4-[4-({[(2-methyl-1,3-benzothiazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-methyl-4-[4-({[(2-methyl-1,3-benzothiazol-5-yl)amino]carbonyl}amino)-3-(trifluoro-methyl)phenoxy]pyridine-2-carboxamide
- N-methyl-4-[3-methyl-4-({[(4-methyl-2-oxo-2H-chromen-7-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- N-methyl-4-[3-methyl-4-({[(2-methyl-1,3-benzothiazol-5yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- 4-[3-fluoro-4-({[(2-methyl-1,3-benzothiazol-5yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-{[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]methyl}-pyridine-2-carboxamide
- 4-{[3-fluoro-4-{{[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]methyl}-N-methylpyridine-2carboxamide
- 4-[2-chloro-4-{{[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide

- N-methyl-4-[3-{{[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxamide
- N-methyl-4-[4-({[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxamide
- 4-[3-({[(2,2-difluoro-1,3-benzodioxol-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-({[(2,2-difluoro-1,3-benzodioxol-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[2-chloro-4-([[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6yl)amino]carbonyl]-amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[2-chloro-4-({[(2,2-difluoro-1,3-benzodioxol-5yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[3-chloro-4-([(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-[3-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-methyl-4-[3-({[(1-methyl-1H-indazol-6yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- 4-(3-{[(2,3-dihydro-1-benzofuran-5-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- N-methyl-4-{3-[([[2-(trifluoromethyl)-1H-benzimidazol-5yl]amino}carbonyl)amino]-phenoxy}pyridine-2-carboxamide
- 4-[4-chloro-3-{{[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-methyl-pyridine-2-carboxamide
- 4-[4-chloro-3-([[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]-carbonyl]amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[4-chloro-3-({[(2,2-difluoro-1,3-benzodioxol-5yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[3-chloro-4-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- 4-[2-chloro-4-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide

- 4-[4-({[(2,2-difluoro-1,3-benzodioxol-5-yl)amino]carbonyl}amino)-3fluorophenoxy]-pyridine-2-carboxamide
- 4-[3-fluoro-4-({[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxamide
- 4-(4-[[(2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- N-methyl-4-[4-({[(1-oxo-2,3-dihydro-1H-inden-5-yl)amino|carbonyl}amino)phenoxyl-pyridine-2-carboxamide
- 5-[3-fluoro-4-({[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxyl-N-methylnicotinamide
- 4-[4-{[(2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino}-3-(trifluoromethyl)phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-[4-({[(1-oxo-2,3-dihydro-1H-inden-5-yl)amino]carbonyl}amino)-3-(trifluoromethyl)phenoxy]pyridine-2-carboxamide
- 4-(3-chloro-4-{[(2,3-dihydro-1H-inden-5ylamino)carbonyl]amino}phenoxy)pyridine-2-carboxamide
- 4-[3-chloro-4-([[(1-oxo-2,3-dihydro-1H-inden-5yl)amino]carbonyl}amino)phenoxy]-pyridine-2-carboxamide
- N-methyl-4-[4-({[(1-methyl-1H-indazol-6yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- 4-(4-{[(1,3-benzothiazol-6-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- N-methyl-4-[4-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- 4-(4-{[(2,3-dihydro-1-benzofuran-5-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- 4-[2,4-dichloro-5-([[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]carbonyl]-amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[2,4-dichloro-5-({(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[3-chloro-4-({[(2,2-difluoro-1,3-benzodioxol-5yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide

- 4-[3-chloro-4-({[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxyl-N-methylpyridine-2-carboxamide;
- 4-(3-chloro-4-[[(2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino)phenoxy)-N-methylpyridine-2-carboxamide
- 4-(3-chloro-4-[[(2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino}phenoxy)-N-methylpyridine-2-carboxamide;
- 4-[3-chloro-4-({[(1-oxo-2,3-dihydro-1H-inden-5yl)amino]carbonyl}amino)phenoxyl-N-methylpyridine-2-carboxamide;
- 4-[2-chloro-4-({[(1-oxo-2,3-dihydro-1H-inden-5-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-(3-chloro-4-{[(2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino}phenoxy)-N-methylpyridine-2-carboxamide
- 4-(3-chloro-4-{[(2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino}phenoxy)-N-methylpyridine-2-carboxamide
- 4-[2,4-dichloro-5-({[(2,2-difluoro-1,3-benzodioxol-5yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-{4-[([[1-(methylsulfonyl)-2,3-dihydro-1H-indol-5yl]amino}carbonyl)amino]-phenoxy}pyridine-2-carboxamide
- N-methyl-4-[3-nitro-4-({[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]-carbonyl}amino)phenoxylpyridine-2-carboxamide
- N-methyl-4-[2-methyl-4-{[[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-6-yl)amino]-carbonyl}amino)phenoxylpyridine-2-carboxamide
- 4-[2,3-difluoro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]-N-methylpyridine-2-carboxamide
- 4-[3,5-difluoro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]-N-methylpyridine-2-carboxamide
- 4-[2,5-difluoro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino]carbonyl}amino)-phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-[4-({[(2,2,3,3-tetrafluoro-2,3-dihydro-1,4-benzodioxin-5yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxamide trifluoroacetate
- 4-[3-fluoro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino]carbonyl}-amino)phenoxy]-pyridine-2-carboxamide

- 4-[3-fluoro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- N-methyl-4-{[5-{{[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino]carbonyl}amino)-quinolin-8-yl]oxy}pyridine-2-carboxamide
- 4-(3-[((1H-indazol-5-ylamino)carbonyl]amino)phenoxy)-N-methylpyridine-2-carboxamide dihydrochloride
- N-[2-(methylamino)-2-oxoethyl]-4-[4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- 4-(3-fluoro-4-{[(quinoxalin-2-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- N-[2-(dimethylamino)-2-oxoethyl]-4-[4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-methyl-4-{3-methyl-4-{{([(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxamide
- Methyl 4-[3-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-7yl)amino]carbonyl}-amino)phenoxyl-pyridine-2-carboxylate
- 4-[3-chloro-4-([[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino]carbonyl]amino)-phenoxy]-N-methylpyridine-2-carboxamide
- 4-[3-chloro-4-{{[[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- 4-(3-{[(1,3-benzodioxol-5-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- N-methyl-4-[3-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino|carbonyl}amino)-phenoxy|pyridine-2-carboxamide
- 4-{3-[[(2,3-dihydro-1,4-benzodioxin-6-ylamino)carbonyl]amino}phenoxy)-N-methylpyridine-2-carboxamide
- 4-[4-chloro-3-([[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino]carbonyl]amino)-phenoxy]-N-methylpyridine-2-carboxamide
- 5-[2-fluoro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]-N-methylnicotinamide
- 4-[2-chloro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide

- 4-[3-chloro-4-([[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- 4-[3-fluoro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- 4-[3-fluoro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- 4-(3-[[(1,3-benzodioxol-5-ylamino)carbonyl]amino}-4-chlorophenoxy)-N-methylpyridine-2-carboxamide
- 4-[4-chloro-3-({[(6-fluoro-4H-1,3-benzodioxin-8yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-(4-{[(1,3-benzodioxol-5-ylamino)carbonyl]amino}-3fluorophenoxy)pyridine-2-carboxamide
- 4-[3-fluoro-4-({[(6-fluoro-4H-1,3-benzodioxin-8-yl)amino]carbonyl}amino)phenoxy]-pyridine-2-carboxamide
- 4-(4-chloro-3-{[(2,3-dihydro-1,4-benzodioxin-6vlamino)carbonyllamino}phenoxy)-N-methylpyridine-2-carboxamide
- 4-[3-{{[(7-fluoro-2,3-dihydro-1,4-benzodioxin-5yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- 4-[3-fluoro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]-N-methylpyridine-2-carboxamide
- 4-(4-{[(1,3-benzodioxol-5-ylamino)carbonyl]amino}phenoxy)-N-methylpyridine-2-carboxamide
- N-methyl-4-[4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- Methyl 4-[4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxylate
- Methyl 5-[4-({((2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino|carbonyl}amino)-phenoxylnicotinate
- 4-[2,4-dichloro-5-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino]carbonyl}amino)-phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-5-[4-([[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]nicotinamide

- 4-(4-[[(1,3-benzodioxol-5-ylamino)carbonyl]amino}-3-chlorophenoxy)-N-methylpyridine-2-carboxamide
- 4-[3-chloro-4-([[(6-fluoro-4H-1,3-benzodioxin-8yl)amino]carbonyl]amino)phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-[2-methyl-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxamide
- N-methyl-4-[3-nitro-4-([[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxamide
- N-methyl-4-[3-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide 1-oxide
- 4-[3-({[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-(2-piperidin-1-ylethyl)pyridine-2-carboxamide
- 4-[3-({[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-(2pyrrolidin-1-ylethyl)pyridine-2-carboxamide
- 4-[3-{{[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-Npvridin-3-vlpvridine-2-carboxamide
- N-[3-(1H-imidazol-1-yl)propyl]-4-[3-({[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)-phenoxy|pyridine-2-carboxamide
- N-(2-piperidin-1-ylethyl)-4-[3-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]-carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-(2-pyrrolidin-1-ylethyl)-4-[3-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]-carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-pyridin-3-yl-4-[3-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxamide
- N-[3-(1H-imidazol-1-yl)propyl]-4-[3-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]-carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-[3-(1H-imidazol-1-yl)propyl]-4-[4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]-carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-(2-pyrrolidin-1-ylethyl)-4-[4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]-carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-(2-piperidin-1-ylethyl)-4-[4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]-carbonyl}amino)phenoxy]pyridine-2-carboxamide

- N-(2-piperazin-1-ylethyl)-4-[4-([[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]-carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-pyridin-2-yl-4-[4-([(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- 4-[4-({[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-(2pyrrolidin-1-ylethyl)pyridine-2-carboxamide
- 4-[4-({[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-(2-piperazin-1-ylethyl)pyridine-2-carboxamide
- 4-[2-methoxy-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- 4-(4-{[(2,3-dihydro-1H-inden-5-ylamino)carbonyl]amino}-2methoxyphenoxy)pyridine-2-carboxamide
- 4-[2,5-difluoro-4-({((2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- 4-[3,5-difluoro-4-({((2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6yl)aminolcarbonyl}amino)-phenoxylpyridine-2-carboxamide
- 4-[3-(aminocarbonyl)-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}-amino)phenoxy]pyridine-2-carboxamide
- N-methyl-4-[3-(methylsulfonyl)-4-({[[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-methyl-4-[3-(methylthio)-4-([[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-6-yl)amino]carbonyl]amino)phenoxy]pyridine-2-carboxamide
- 4-[3-fluoro-4-([[(6-nitro-1,3-benzothiazol-2yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-[4-({[(6-nitro-1,3-benzothiazol-2yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- 4-[4-([[(4,6-difluoro-1,3-benzothiazol-2-yl)amino]carbonyl]amino)phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-[4-({[(2-methyl-1,3-benzoxazol-6yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- 4-(4-[[(2,3-dihydro-1H-inden-4-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide

- 4-[4-({[(2,2-difluoro-1,3-benzodioxol-4-yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- N-methyl-4-[4-({[(2-methyl-2H-indazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- 4-(4-[[({1-[2-(diethylamino)ethyl]-1H-indazol-5-yl}amino)carbonyl]amino}-3fluorophenoxy)-N-methylpyridine-2-carboxamide
- N-methyl-4-[4-({[(2-methyl-1H-indol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-{4-[(2-acetylpyridin-4-yl)oxy]phenyl}-N'-(1-methyl-1H-indazol-5-yl)urea
- N-[2-(dimethylamino)-2-oxoethyl]-4-[4-({[(1-methyl-1H-indazol-5-yl)amino]carbonyl}-aminophenoxy]pyridine-2-carboxamide
- N-methyl-4-[4-({[(2-methyl-1,3-benzothiazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-methyl-4-{[4-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]methyl}-pyridine-2-carboxamide
- 4-(3-[[(1H-1,2,3-benzotriazol-5-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- Methyl 4-[3-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxylate
- 4-(4-{[(1H-1,2,3-benzotriazol-5-ylamino)carbonyl]amino}phenoxy)-Nmethylpyridine-2-carboxamide
- 4-(4-[[(1H-indazol-6-ylamino)carbonyl]amino}phenoxy)-N-methylpyridine-2-carboxamide
- N-methyl-4-{4-[([2-(trifluoromethyl)-1H-benzimidazol-5yl]amino}carbonyl)amino]-phenoxy}pyridine-2-carboxamide
- 4-[4-{{[(1-ethyl-2-methyl-1H-benzimidazol-5yl)amino]carbonyl}amino)phenoxy]-N-methylpyridine-2-carboxamide
- Methyl 4-[4-({[(1-methyl-1H-indazol-5yl)amino|carbonyl}amino)phenoxy]pyridine-2-carboxylate
- 4-[2-chloro-4-({[(2,2,4,4-tetrafluoro-4H-1,3-benzodioxin-7yl)amino]carbonyl}amino)-phenoxy]-N-methylpyridine-2-carboxamide
- 4-(4-{[(2,3-dihydro-1,4-benzodioxin-6-ylamino)carbonyl]amino}phenoxy)-N-[3-(1H-imidazol-1-yl)propyl]pyridine-2-carboxamide

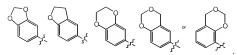
- 4-(4-{[(2,3-dihydro-1,4-benzodioxin-6-ylamino)carbonyl]amino}phenoxy)-N-(2-pyrrolidin-1-ylethyl)pyridine-2-carboxamide
- N-[3-(1H-imidazol-1-yl)propyl]-4-[4-({[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)-phenoxy]pyridine-2-carboxamide
- 4-[4-({[(1-methyl-1H-indazol-5-yl)amino]carbonyl}amino)phenoxy]-N-(2-piperidin-1-ylethyl)pyridine-2-carboxamide
- N-cyclopropyl-4-[4-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide
- N-(cyclopropylmethyl)-4-[4-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]-pyridine-2-carboxamide
- N-cyclobutyl-4-[4-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]pyridine-2-carboxamide or
- Methyl-N-({4-[4-({[(1-methyl-1H-indazol-5yl)amino]carbonyl}amino)phenoxy]pyridin-2-yl}carbonyl)glycinate_
- 13) (original) A pharmaceutical composition which comprises an effective amount of at least one compound of claim1 and a physiologically acceptable carrier.
 - 14) (cancelled)
 - 15) (cancelled)
 - (cancelled)
 - 17) (cancelled)
 - 18) (cancelled)
 - 19) (cancelled)
 - 20) (cancelled)
 - 21) (cancelled)
 - 22) (cancelled)
 - 23) (cancelled)
 - 24) (previously presented) A compound of formula (I):

wherein

Q is C(O)R⁴, C(O)OR⁴ or C(O)NR⁴R⁵;

wherein A is a bicyclic heterocycle which is:

- (1) benzimidazol-5-yl
- (2) benzimidazol-6-yl
- (3) 1,3-benzothiazol-2-yl
- (4) 1,3-benzothiazol-5-yl
- (5) 1,3-benzothiazol-6-yl
- (6) 1,2,3-benzotriazol-5-vl
- (7) 1,3-benzoxazol-2-yl
- (8) 1,3-benzoxazol-6-yl
- (9) 2,3-dihydro-1H-indol-5-yl
- (10) 2,3-dihydro-1H-indol-6-yl
- (11) 2,3-dihydro-1H-inden-4-yl
- (12) 2,3-dihydro-1H-inden-5-yl
- (13) 1,1-dioxido-2,3-dihydro-1-benzothien-6-yl
- (14) 1H-indazol-5-yl
- (15) 2H-indazol-5-yl
- (16) 1H-indazol-6-yl
- (17) 1H-indol-5-yl
- (18) 2-oxo-2H-chromen-7-yl
- (19) 1-oxo-2,3-dihydro-1H-inden-5-yl
- (20) quinoxalin-2-yl
- (21) quinoxalin-6-yl, or
- (22) a group of the formula



optionally substituted with 1-4 substituents which are independently R^1 , OR^1 , $S(O)_pR^1$, $C(O)R^1$, $C(O)OR^1$, $C(O)NR^1R^2$, halogen, oxo, cyano, or nitro

B is phenyl, optionally substituted with 1-4 substituents which are independently C_1 - C_5 linear or branched alkyl, C_1 - C_5 linear or branched haloalkyl, C_1 - C_3 alkoxy, hydroxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, carboxyamide, halogen, cyano, nitro or $S(O)_oR^7$;

Lis:

- (a) -(CH₂)_m-O-(CH₂)_I-,
- (b) $-(CH_2)_m-(CH_2)_{i-}$,
- (c) -(CH₂)_m-C(O)-(CH₂)_I-,
- (d) -(CH₂)_m-NR³-(CH₂)_I-,
- (e) $-(CH_2)_{m}$ $NR^3C(O)-(CH_2)_{l}$ -,
- $(f) (CH_2)_m S (CH_2)_{l^-},$
- (g) -(CH₂)_m-C(O)NR³ -(CH₂)_I-, or
- (h) a single bond;

m and I are integers independently selected from 0-4;

M is a pyridine ring, optionally substituted with 1-3 substituents which are independently C_1 - C_5 linear or branched alkyl, C_1 - C_5 linear or branched haloalkyl, C_1 - C_3 alkoxy, hydroxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, halogen, or nitro;

Q is C(O)R⁴, C(O)OR⁴ or C(O)NR⁴R⁵;

each of R1, R2, R3, R4 and R5, is independently:

- (a) hydrogen,
- (b) C₁-C₅ linear, branched, or cyclic alkyl,
- (c) phenyl,
- (d) C₁-C₃ alkyl-phenyl,
- (e) up to per-halo substituted C₁-C₅ linear or branched alkyl,
- (f) -(CH₂)_q·X, wherein X is a 5 or 6 membered heterocyclic ring, containing at least one atom selected from oxygen, nitrogen and sulfur, which is saturated, partially saturated, or aromatic, or a 8-10 membered bicyclic heteroaryl having 1-4 heteroatoms which are O, N or S, or
- (g) -(CH₂)_q-Y, where Y is C(O)R⁶, C(O)OR⁶ and C(O)NR⁶R⁷;

each of R⁶ - R⁷ is independently :

- (a) hydrogen,
- (b) C₁-C₅ linear, branched, or cyclic alkyl,
- (c) phenyl,
- (d) C1-C3 alkyl-phenyl, or
- (e) up to per-halo substituted C₁-C₅ linear or branched alkyl;

each of R^1 , R^2 , R^3 , R^4 , R^5 , R^6 and R^7 , other than per-halo substituted C_1 - C_5 linear or branched alkyl, is optionally substituted with 1-3 substituents which are independently C_1 - C_5 linear or branched alkyl, up to perhalo substituted C_1 - C_5 linear or branched alkyl, C_1 - C_3 alkoxy, hydroxy, carboxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, halogen, cyano, or nitro;

p is an integer selected from 0, 1, or 2; and

g is an integer selected from 1, 2, 3, or 4,

or a pharmaceutically acceptable salt of formula I or an oxidized derivative of formula I wherein one or more urea nitrogens are substituted with a hydroxyl group, or an oxidized derivative of formula I wherein the nitrogen atom of

pyridine ring M is in the oxide form, or a methyl, ethyl, propyl, isopropyl, butyl, isobutyl, pentyl ester or phenyl C₁-C₅ alkyl ester of formula I at a carboxylic acid group or amide group.

25) (original) A compound of claim 24 wherein A is selected from

- (1) benzimidazol-5-yl
- (2) benzimidazol-6-yl
- (8) 1.3-benzoxazol-6-vl
- (9) 2,3-dihydro-1H-indol-5-yl
- (10) 2,3-dihydro-1H-indol-6-yl
- (11) 2,3-dihydro-1H-inden-4-yl
- (12) 2,3-dihydro-1H-inden-5-yl
- (13) 1,1-dioxido-2,3-dihydro-1-benzothien-6-vl
- (14) 1H-indazol-5-yl
- (15) 2H-indazol-5-yl
- (16) 1H-indazol-6-yl
- (17) 1H-indol-5-yl
- (18) quinoxalin-2-yl
- (19) quinoxalin-6-yl, and
- (20) a group of the formula

26) (original) A compound of claim 24 wherein the optional substituents on bicyclic heterocycle A are independently R^1 , OR^1 , and halogen.

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- 27) (previously presented) A compound as in claim 26 wherein B is phenyl optionally substituted with 1-4 substituents which are halogen.
 - 28) (original) A compound of claim 27 wherein L is -O-.
- 29) (original) A compound of claim 28 wherein Q is $C(O)NR^4R^5$ and each of R^4 and R^5 is independently hydrogen or C_1 - C_5 alkyl.
 - 30) (previously presented) A compound of formula (I):

wherein

A is a bicyclic heterocycle which is:

- (1) benzimidazol-5-yl
- (2) benzimidazol-6-yl
- (8) 1,3-benzoxazol-6-yl
- (9) 2,3-dihydro-1H-indol-5-yl
- (10) 2,3-dihydro-1H-indol-6-yl
- (11) 2,3-dihydro-1H-inden-4-yl
- (12) 2,3-dihydro-1H-inden-5-yl
- (13) 1,1-dioxido-2,3-dihydro-1-benzothien-6-yl
- (14) 1H-indazol-5-yl
- (15) 2H-indazol-5-yl
- (16) 1H-indazol-6-yl
- (17) 1H-indol-5-yl
- (18) quinoxalin-2-yl
- (19) quinoxalin-6-yl, and
- (20) a group of the formula

optionally substituted with 1-4 substituents which are independently R^1 , OR^1 , $S(O)_pR^1$, $C(O)R^1$, $C(O)OR^1$, $C(O)NR^1R^2$, halogen, oxo, cyano, or nitro B is phenyl, optionally substituted with halogen,

L is -O-,

M is a pyridine ring substituted only with Q, Q is C(O)NHR⁵ and R⁵ is independently hydrogen or C₁-C₅ alkyl, and p is an integer selected from 0, 1, or 2

or a pharmaceutically acceptable salt of formula I or an oxidized derivative of formula I wherein one or more urea nitrogens are substituted with a hydroxyl group, or an oxidized derivative of formula I wherein the nitrogen atom of pyridine ring M is in the oxide form, or a methyl, ethyl, propyl, isopropyl, butyl, isobutyl, pentyl ester or phenyl C₁-C₅ alkyl ester of formula I at a carboxylic acid group or amide group.

31) (cancelled)

32) (New) A compound of formula (I):

wherein

A is a bicyclic heterocycle which is:

- (1) benzimidazolyl
- (2) 1,3-benzothiazolyl
- (3) 1.2.3-benzotriazolvl

- (4) 1,3-benzoxazolyl
- (5) 2,3-dihydro-1H-indolyl
- (6) 2,3-dihydro-1H-indenyl
- (7) 1.1-dioxido-2.3-dihydro-1-benzothienvl
- (8) 1H-indazolyl
- (9) 2H-indazolyl
- (10) 1H-indolyl
- (11) 2H-chromenyl
- (12) quinoxalinyl or
- (13) a group of the formula

optionally substituted with 1-4 substituents which are independently R^1 , OR^1 , $S(O)_nR^1$, $C(O)R^1$, $C(O)OR^1$, $C(O)NR^1R^2$, halogen, oxo, cyano, or nitro;

B is quinolinyl, optionally substituted with 1-4 substituents which are independently C_1 - C_5 linear or branched alkyl, C_1 - C_5 linear or branched haloalkyl, C_1 - C_3 alkoxy, hydroxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, carboxyamide, halogen, cyano, nitro or $S(O)_oR^7$;

Lis:

- (a) -(CH₂)_m-O-(CH₂)_I-,
- (b) -(CH₂)_m-(CH₂)_I-,
- (c) $-(CH_2)_m C(O) (CH_2)_{1-}$,
- (d) -(CH₂)_m-NR³-(CH₂)_I-,
- (e) -(CH₂)_m- NR³C(O)-(CH₂)_I-,
- (f) $-(CH_2)_m-S-(CH_2)_{l^-}$,
- (g) -(CH₂)_m-C(O)NR³ -(CH₂)_i-, or
- (h) a single bond;

m and I are integers independently selected from 0-4:

M is a pyridine ring, optionally substituted with 1-3 substituents which are independently C_1 - C_5 linear or branched alkyl, C_1 - C_5 linear or branched haloalkyl, C_1 - C_3 alkoxy, hydroxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, halogen, or nitro;

Q is $C(O)R^4$, $C(O)OR^4$ or $C(O)NR^4R^{5.}$;

each of R1, R2, R3, R4 and R5 is independently:

- (a) hydrogen,
- (b) C₁-C₅ linear, branched, or cyclic alkyl.
- (c) phenyl,
- (d) C₁-C₃ alkyl-phenyl,
- (e) up to per-halo substituted C₁-C₅ linear or branched alkyl,
- (f) -(CH₂)_q·X, wherein X is a 5 or 6 membered heterocyclic ring, containing at least one atom selected from oxygen, nitrogen and sulfur, which is saturated, partially saturated, or aromatic, or a 8-10 membered bicyclic heteroaryl having 1-4 heteroatoms which are O. N or S. or
- (g) -(CH₂)_{σ}-Y, where Y is C(O)R⁶, C(O)OR⁶ and C(O)NR⁶R⁷:

each of R⁶ - R⁷ is independently:

- (a) hydrogen,
- (b) C₁-C₅ linear, branched, or cyclic alkyl,
- (c) phenyl,
- (d) $C_1\text{-}C_3$ alkyl-phenyl, or
- (e) up to per-halo substituted C₁-C₅ linear or branched alkyl;

each of R^1 , R^2 , R^3 , R^4 , R^5 , R^6 and R^7 , other than per–halo substituted C_1 - C_5 linear or branched alkyl, is optionally substituted with 1-3 substitutents which are independently C_1 - C_5 linear or branched alkyl, up to perhalo substituted C_1 -

C₅ linear or branched alkyl, C₁-C₃ alkoxy, hydroxy, carboxy, amino, C₁-C₃ alkylamino, C₁-C₆ dialkylamino, halogen, cyano, or nitro;

p is an integer selected from 0, 1, or 2; and

q is an integer selected from 1, 2, 3, or 4

or a pharmaceutically acceptable salt of formula I or an oxidized derivative of formula I wherein one or more urea nitrogens are substituted with a hydroxyl group, or an oxidized derivative of formula I wherein the nitrogen atom of pyridine ring M is in the oxide form, or a methyl, ethyl, propyl, isopropyl, butyl, isobutyl, pentyl ester or phenyl C₁-C₅ alkyl ester of formula I at a carboxylic acid group or amide group.

33) (New) A compound of claim 1 wherein A and B follow one of the following combinations:

A= 1H-benzimidazol-5-yl; and B= quinolinyl,

A= 1H-benzimidazol-6-yl; and B= quinolinyl,

A= 1,3-benzodioxin-6-yl; and B= quinolinyl,

A= 1,3-benzodioxin-7-yl; and B= quinolinyl,

A= 1,3-benzodioxin-8-yl; and B= quinolinyl,

A= 1,3-benzodioxol-4-yl; and B= quinolinyl,

A= 1,3-benzodioxol-5-yl; and B= quinolinyl,

A= 1,3-benzothiazol-2-yl; and B= quinolinyl,

A= 1,3-benzothiazol-5-yl; and B= quinolinyl, A= 1,3-benzothiazol-6-yl; and B= quinolinyl,

A= 1,2,3-benzotriazol-5-yl; and B= quinolinyl,

A= 1,3-benzoxazol-2-yl; and B= quinolinyl, or

A= 1,3-benzoxazol-6-yl; and B= quinolinyl.

34) (New) A compound of claim 1 wherein A and B follow one of the following combinations:

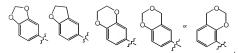
- A= 2,3-dihydro-1,4-benzodioxin-5-yl; and B= quinolinyl,
- A= 2,3-dihydro-1,4-benzodioxin-6-yl; and B= quinolinyl,
- A= 2,3-dihydro-1-benzofuran-5-yl; and B= quinolinyl,
- A= 2,3-dihydro-1H-indol-5-yl; and B= quinolinyl,
- A= 2,3-dihydro-1H-indol-6-yl; and B= quinolinyl,
- A= 2,3-dihydro-1H-inden-4-yl; and B= quinolinyl,
- A= 2,3-dihydro-1H-inden-5-yl; and B= quinolinyl,
- A= 1,1-dioxido-2,3-dihydro-1-benzothien-6-yl; and B= quinolinyl.
- 35) (New) A compound of claim 1 wherein A and B follow one of the following combinations:
 - A= 1H-indazol-5-yl; and B= quinolinyl,
 - A= 2H-indazol-5-yl; and B= quinolinyl,
 - A= 1H-indazol-6-yl; and B= quinolinyl,
 - A= 1H-indol-5-yl; and B= quinolinyl,
 - A= 2-oxo-2H-chromen-7-yl; and B= quinolinyl or
 - A= 1-oxo-2,3-dihydro-1H-inden-5-yl and B=quinolinyl.
- 36) (New) A compound of claim 1 wherein A and B follow one of the following combinations:
 - A= quinoxalin-2-yl; and B= quinolinyl or
 - A= quinoxalin-6-yl; and B= quinolinyl.
 - 37) (New) A compound as in claim 32 wherein L is -O- or -S-.
- 38) (New) A pharmaceutical composition which comprises an effective amount of at least one compound of claim 32 and a physiologically acceptable carrier.
 - 39) (New) A compound of formula (I):

wherein

Q is C(O)R⁴, C(O)OR⁴ or C(O)NR⁴R⁵;

wherein A is a bicyclic heterocycle which is:

- (1) benzimidazol-5-yl
- (2) benzimidazol-6-yl
- (3) 1,3-benzothiazol-2-yl
- (4) 1,3-benzothiazol-5-yl
- (5) 1,3-benzothiazol-6-yl
- (6) 1,2,3-benzotriazol-5-vl
- (7) 1,3-benzoxazol-2-yl
- (8) 1,3-benzoxazol-6-yl
- (9) 2,3-dihydro-1H-indol-5-yl
- (10) 2,3-dihydro-1H-indol-6-yl
- (11) 2,3-dihydro-1H-inden-4-yl
- (12) 2,3-dihydro-1H-inden-5-yl
- (13) 1,1-dioxido-2,3-dihydro-1-benzothien-6-yl
- (14) 1H-indazol-5-yl
- (15) 2H-indazol-5-yl
- (16) 1H-indazol-6-yl
- (17) 1H-indol-5-yl
- (18) 2-oxo-2H-chromen-7-yl
- (19) 1-oxo-2,3-dihydro-1H-inden-5-yl
- (20) quinoxalin-2-yl
- (21) quinoxalin-6-yl, or
- (22) a group of the formula



optionally substituted with 1-4 substituents which are independently R^1 , OR^1 , $S(O)_pR^1$, $C(O)R^1$, $C(O)OR^1$, $C(O)NR^1R^2$, halogen, oxo, cyano, or nitro

B is quinolinyl, optionally substituted with 1-4 substituents which are independently C_1 - C_5 linear or branched alkyl, C_1 - C_5 linear or branched haloalkyl, C_1 - C_3 alkoxy, hydroxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, carboxyamide, halogen, cyano, nitro or $S(O)_oR^7$;

L is:

- (a) -(CH₂)_m-O-(CH₂)_I-,
- (b) $-(CH_2)_m-(CH_2)_{l}-$,
- (c) -(CH₂)_m-C(O)-(CH₂)_I-,
- (d) -(CH₂)_m-NR³-(CH₂)_I-,
- (e) -(CH₂)_m- NR³C(O)-(CH₂)_I-,
- $(f) \ \hbox{-} (CH_2)_m \hbox{-} S \hbox{-} (CH_2)_{l^-},$
- (g) $-(CH_2)_m-C(O)NR^3-(CH_2)_{I^-}$, or
- (h) a single bond;

m and I are integers independently selected from 0-4;

M is a pyridine ring, optionally substituted with 1-3 substituents which are independently C_1 - C_5 linear or branched alkyl, C_1 - C_5 linear or branched haloalkyl, C_1 - C_3 alkoxy, hydroxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, halogen, or nitro;

Q is $C(O)R^4$, $C(O)OR^4$ or $C(O)NR^4R^5$;

each of R1, R2, R3, R4 and R5, is independently:

- (a) hydrogen,
- (b) C₁-C₅ linear, branched, or cyclic alkyl,
- (c) phenyl,
- (d) C₁-C₃ alkyl-phenyl,
- (e) up to per-halo substituted C₁-C₅ linear or branched alkyl,
- (f) -(CH₂)_q·X, wherein X is a 5 or 6 membered heterocyclic ring, containing at least one atom selected from oxygen, nitrogen and sulfur, which is saturated, partially saturated, or aromatic, or a 8-10 membered bicyclic heteroaryl having 1-4 heteroatoms which are O, N or S, or
- (g) -(CH₂)_q-Y, where Y is C(O)R⁶, C(O)OR⁶ and C(O)NR⁶R⁷;

each of R⁶ - R⁷ is independently :

- (a) hydrogen,
- (b) C₁-C₅ linear, branched, or cyclic alkyl,
- (c) phenyl,
- (d) C1-C3 alkyl-phenyl, or
- (e) up to per-halo substituted C₁-C₅ linear or branched alkyl;

each of R^1 , R^2 , R^3 , R^4 , R^5 , R^6 and R^7 , other than per-halo substituted C_1 - C_5 linear or branched alkyl, is optionally substituted with 1-3 substituents which are independently C_1 - C_5 linear or branched alkyl, up to perhalo substituted C_1 - C_5 linear or branched alkyl, C_1 - C_3 alkoxy, hydroxy, carboxy, amino, C_1 - C_3 alkylamino, C_1 - C_6 dialkylamino, halogen, cyano, or nitro;

p is an integer selected from 0, 1, or 2; and

q is an integer selected from 1, 2, 3, or 4,

or a pharmaceutically acceptable salt of formula I or an oxidized derivative of formula I wherein one or more urea nitrogens are substituted with a hydroxyl group, or an oxidized derivative of formula I wherein the nitrogen atom of pyridine ring M is in the oxide form, or a methyl, ethyl, propyl, isopropyl, butyl, isobutyl, pentyl ester or phenyl C₁-C₅ alkyl ester of formula I at a carboxylic acid group or amide group.

40) (New) A compound of claim 39 wherein A is selected from

- (1) benzimidazol-5-yl
- (2) benzimidazol-6-yl
- (8) 1,3-benzoxazol-6-yl
- (9) 2,3-dihydro-1H-indol-5-yl
- (10) 2,3-dihydro-1H-indol-6-yl
- (11) 2,3-dihydro-1H-inden-4-yl
- (12) 2,3-dihydro-1H-inden-5-yl
- (13) 1,1-dioxido-2,3-dihydro-1-benzothien-6-vl
- (14) 1H-indazol-5-yl
- (15) 2H-indazol-5-yl
- (16) 1H-indazol-6-yl
- (17) 1H-indol-5-yl
- (18) quinoxalin-2-vl
- (19) quinoxalin-6-yl, and
- (20) a group of the formula

- 41) (New) A compound of claim 39 wherein the optional substituents on bicyclic heterocycle A are independently R¹, OR¹, and halogen.
 - 42) (New) A compound of claim 41 wherein L is -O-.

 $\mbox{43)} \quad \mbox{(New) A compound of claim 42 wherein Q is } \mbox{ C(O)NR}^4 R^5 \mbox{ and} \\ \mbox{each of R}^4 \mbox{ and R}^5 \mbox{ is independently hydrogen or C}_1 - \mbox{C}_5 \mbox{ alkyl}.$